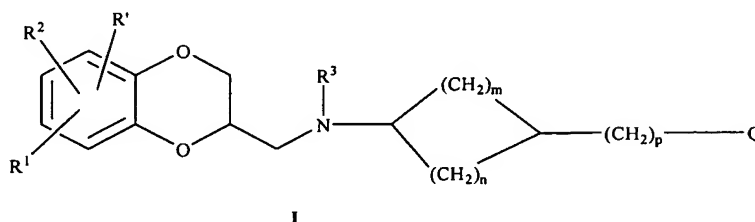


This listing of claims will replace all prior versions, and listings, of claims in the application.

**Listing of Claims**

1. (*currently amended*) A compound of Formula I:



wherein

$R^2$ ,  $R^1$  and  $R^2$  are, independently, hydrogen, halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon atoms, or alkanesulfonamido of 1 to 6 carbon atoms; or  $R^1$  and  $R^2$ , taken together, form methylenedioxy, ethylenedioxy or propylenedioxy;

**$R^1$  is halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon atoms, or alkanesulfonamido of 1 to 6 carbon atoms;**

$R^3$  is hydrogen or alkyl of 1 to 6 carbon atoms;

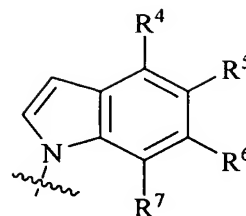
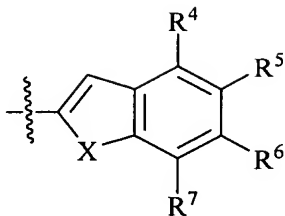
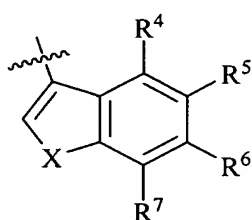
m is 1 to 3;

n is 1 or 2;

p is 0 to 3

with the proviso that when p is 0, both m and n may not be 2;

Q is a heteroaryl moiety chosen from:



$R^4$ ,  $R^5$ ,  $R^6$  and  $R^7$  are independently selected from hydrogen, halo, cyano, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms;

X is  $NR^8$ , ~~O~~ or ~~S~~; and

$R^8$  is hydrogen or alkyl of 1 to 6 carbon atoms;  
or a pharmaceutically acceptable salt thereof.

2. *(original)* A compound according to claim 1, wherein  $R^1$  and  $R^2$  are, independently, hydrogen, halo, cyano, carboxamido, trifluoromethyl, amino, alkyl of one to six carbon atoms or alkoxy of one to six carbon atoms.
3. *(original)* A compound according to claim 1, wherein  $R^1$  is alkoxy of one to six carbon atoms and is attached to position 8 of the benzodioxan moiety.
4. *(original)* A compound according to claim 1, wherein  $R^2$  is hydrogen.
5. *(original)* A compound according to claim 1, wherein  $R^4$ ,  $R^5$ ,  $R^6$  and  $R^7$  are independently hydrogen, halo, cyano, alkyl of one to six carbon atoms, or alkoxy of one to six carbon atoms.
6. *(original)* A compound according to claim 1, wherein  $R^4$ ,  $R^5$ ,  $R^6$  and  $R^7$  are independently hydrogen, halo or cyano.
7. *(original)* A compound according to claim 1, wherein m and n are, independently 1 or 2.
8. *(original)* A compound according to claim 1, wherein m is 1 and n is 2.

9. *(original)* A compound according to claim 1, wherein p is 0 or 1.
10. *(original)* A compound according to claim 1, wherein p is 0.
11. *(currently amended)* A compound according to claim 1, wherein ~~X is NR<sup>8</sup> and~~ R<sup>8</sup> hydrogen or alkyl of 1 to 3 carbons.
12. *(original)* A compound according to claim 1, wherein R<sup>3</sup> is hydrogen or alkyl of 1 to 3 carbons.
13. *(previously presented)* A compound according to claim 1, wherein said compound is N-[(cis)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
14. *(original)* A compound according to claim 1, wherein said compound is N-[(1R,3S)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
15. *(original)* A compound according to claim 1, wherein said compound is N-[(1S,3R)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
16. *(original)* A compound according to claim 1, wherein said compound is N-[(trans)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
17. *(original)* A compound according to claim 1, wherein said compound is N-[(1S,3S)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.

18. (*original*) A compound according to claim 1, wherein said compound is N-[(1R,3R)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
19. (*original*) A compound according to claim 1, wherein said compound is N-[(cis)-3-(5-fluoro-1-methyl-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
20. (*original*) A compound according to claim 1, wherein said compound is N-[(trans)-3-(5-fluoro-1-methyl-1H-indol-3-yl)cyclopentyl]-N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amine or a pharmaceutically acceptable salt thereof.
21. (*original*) A compound according to claim 1, wherein said compound is 3-[(cis)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.
22. (*original*) A compound according to claim 1, wherein said compound is 3-[(trans)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.
23. (*original*) A compound according to claim 1, wherein said compound is 3-[(cis)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1-methyl-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.
24. (*original*) A compound according to claim 1, wherein said compound is 3-[(trans)-3-({[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}amino)cyclopentyl]-1-methyl-1H-indole-5-carbonitrile or a pharmaceutically acceptable salt thereof.
25. (*original*) A compound according to claim 1, wherein said compound is N-{[(2S)-8-methoxy-2,3-dihydro-1,4-benzodioxin-2-yl]methyl}-N-[(cis)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]amine or a pharmaceutically acceptable salt thereof.

26. (*original*) A compound according to claim 1, wherein said compound is N-{[(2S)-8-ethoxy-2,3-dihydro-1,4-benzo-dioxin-2-yl]methyl}-N-[(trans)-3-(5-fluoro-1H-indol-3-yl)cyclopentyl]amine or a pharmaceutically acceptable salt thereof.

27. (*cancelled*)

28. (*cancelled*)

29. (*cancelled*)

30. (*cancelled*)

31. (*original*) A pharmaceutical composition, comprising:  
an effective amount of a compound according to claim 1 or a pharmaceutically acceptable salt thereof; and  
a pharmaceutically acceptable carrier or excipient.